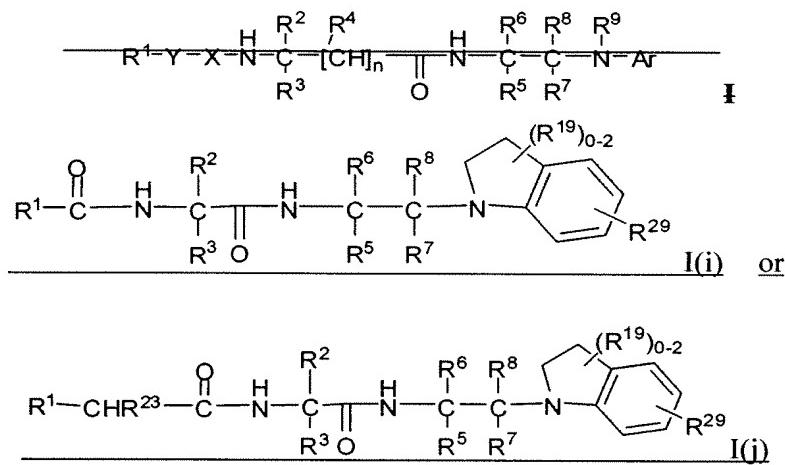


CLAIMS

1. (currently amended) A compound of Formula I(i) or I(j):



or a pharmaceutically acceptable salt thereof, wherein:

R^1 is $\text{C}_6\text{-C}_{10}$ aryl substituted with 0-3 R^{1a} , or a $\text{C}_3\text{-C}_8$ cycloalkyl substituted with 0-2 R^{1b} , wherein said $\text{C}_3\text{-C}_8$ cycloalkyl is saturated or unsaturated;

each R^{1a} is independently a member selected from the group consisting of H, $\text{C}_1\text{-C}_3$ perfluoroalkyl, $\text{C}_3\text{-C}_7$ cycloalkyl, F, Cl, Br, CN, NO_2 , OR^{10} , SCH_3 , $\text{S}(=\text{O})\text{CH}_3$, $\text{S}(=\text{O})_2\text{R}^{10}$, $\text{NR}^{11}\text{R}^{12}$, acetyl, $\text{C}(=\text{O})\text{OR}^{13}$, $\text{C}(=\text{O})\text{NR}^{13}\text{R}^{14}$, $\text{S}(=\text{O})_2\text{NR}^{13}\text{R}^{14}$, phenyl substituted with 0-3 R^{15} , and a $\text{C}_1\text{-C}_4$ alkyl substituted with 0-2 R^{16} ;

each R^{1b} is independently a member selected from the group consisting of H, OH, F, Cl, acetyl, $=\text{O}$, $\text{C}_1\text{-C}_6$ alkyl, $\text{C}_1\text{-C}_6$ alkoxy, CF_3 and OCF_3 ;

R^2 is a member selected from the group consisting of a phenyl substituted with 0-3 R^{15} , a $\text{C}_1\text{-C}_6$ alkyl substituted with 0-2 R^{2a} , ~~wherein said C₁-C₆ alkyl optionally contains a heteroatom selected from the group consisting of O, S, and S(=O)₂~~, a $\text{C}_2\text{-C}_6$ alkenyl, a $\text{C}_2\text{-C}_6$ alkynyl, a $\text{C}_3\text{-C}_7$ cycloalkyl substituted with 0-2 R^{19} , and a $\text{C}_7\text{-C}_{11}$ bicycloalkyl substituted with 0-2 R^{19} ;

each R^{2a} is independently a member selected from the group consisting of a $\text{C}_6\text{-C}_{10}$ aryl substituted with 0-3 R^{15} , a $\text{C}_3\text{-C}_8$ cycloalkyl substituted with 0-2 R^{19} , and a $\text{C}_7\text{-C}_{11}$ bicycloalkyl substituted with 0-2 R^{19} ;

R^3 is a member selected from the group consisting of H and $\text{C}_1\text{-C}_4$ alkyl;

~~subscript n is 0 or 1;~~

~~R⁴ is a member selected from the group consisting of H and C₁-C₆ alkyl;~~

~~R⁵ is a member selected from the group consisting of H, C₃-C₇ cycloalkyl, C₂-C₆ alkenyl, C₂-C₆ alkyne, phenyl substituted with 0-2 R¹⁵; and a C₁-C₆ alkyl substituted with 0-2 R¹⁸, wherein said C₁-C₆ alkyl optionally contains a heteroatom selected from the group consisting of O, S, S(=O), S(=O)₂ and NR¹⁷;~~

~~Y is a member independently selected from the group consisting of a bond and (CR²⁰R²¹)_m-W-(CR²²R²³)_p;~~

~~subscript p is 1 or 2;~~

~~subscript m is 0 or 1;~~

~~W is a member independently selected from the group consisting of a bond, O, S, -S(=O), S(=O)₂ and NR¹²;~~

~~X is selected from the group consisting of C(=O), OC(=O), NR²⁴C(=O) and -S(=O)₂;~~

each of R⁶, R⁷, ~~and R⁸ and R⁹~~ is independently a member selected from the group consisting of H and C₁-C₄ alkyl;

~~Ar is a member selected from the group consisting of phenyl substituted with 0-3 R²⁹, and 5-to-6-membered heteroaryl containing 1-to-4 heteroatoms each independently a member selected from the group consisting of N, O and S, wherein said heteroaryl is substituted with 0-3 R²⁹;~~

each R¹⁰ is independently a member selected from the group consisting of H, C₃-C₇ cycloalkyl, a C₁-C₃ perfluoroalkyl, a C₁-C₄ alkyl substituted with 0-1 R²⁵, and a phenyl substituted with 0-3 R¹⁵;

each R¹¹ is independently a member selected from the group consisting of H, 'BOC, Cbz, C₃-C₈ cycloalkyl, (C₁-C₆ alkyl)-C(=O)-, (C₁-C₆ alkyl)-S(=O)₂- and a C₁-C₆ alkyl;

each of R¹², R¹³ and R¹⁴ is independently a member selected from the group consisting of H and C₁-C₄ alkyl;

each R¹⁵ is independently a member selected from the group consisting of H, OH, F, Cl, Br, I, CN, NO₂, COOR¹³, C(=O)NR¹³R¹⁴, S(=O)₂NR¹³R¹⁴, acetyl, -SCH₃, -S(=O)CH₃, -S(=O)₂CH₃, NR²⁶R²⁷, C₁-C₆ alkoxy, C₁-C₃ perfluoroalkyl, C₁-C₃ perfluoroalkoxy and a C₁-C₆ alkyl;

each R¹⁶ is independently a member selected from the group consisting of H, OH, COOR¹³, C(=O)NR¹³R¹⁴, S(=O)₂NR¹³R¹⁴, acetyl, -SCH₃, -S(=O)CH₃, -S(=O)₂CH₃, C₁-C₆ alkoxy, NR²⁶R²⁷, and a phenyl substituted with 0-3 R¹⁵;

~~R¹⁷ is a member selected from the group consisting of H and C₁-C₄ alkyl;~~

each R¹⁸ is independently a member selected from the group consisting of H, OH, F, Cl, CN, NO₂, C(=O)OR³⁰, C(=O)NR¹³R¹⁴, NR¹¹R¹², a C₁-C₃ perfluoroalkyl, a C₁-C₃ perfluoroalkoxy, a phenyl substituted with 0-3 R¹⁵; and C₃-C₈ cycloalkyl;

each R¹⁹ is [[a]] independently a member selected from the group consisting of C₁-C₄ alkyl, F, Cl, and C₁-C₄ alkoxy, CF₃ and OCF₃;

~~each of R²⁰, R²¹, R²² and R²³ is independently a member selected from the group consisting of a bond, H, F, OH, C₁-C₄ alkyl, and C₁-C₃ alkylhydroxy;~~

~~R²⁴ is a member selected from the group consisting of H and C₁-C₄ alkyl;~~

each R²⁵ is independently a member selected from the group consisting of H, C₃-C₇ cycloalkyl, and a phenyl substituted with 0-3 R¹⁵;

each R²⁶ is independently a member selected from the group consisting of H, C₁-C₄ alkyl, (C₁-C₄ alkyl)-C(=O)- and (C₁-C₄ alkyl)-S(=O)₂-;

each R²⁷ is independently a member selected from the group consisting of H and C₁-C₄ alkyl;

each R²⁸ is independently a member selected from the group consisting of H, a C₁-C₆ alkyl, C₃-C₈ cycloalkyl, a phenyl substituted with 0-3 R¹⁵, and a benzyl substituted with 0-2 R¹⁵;

each R²⁹ is independently a member selected from the group consisting of H, F, Cl, Br, I, CN, NO₂, OR²⁸, SR²⁸, S(=O)R²⁸, S(=O)₂R²⁸, S(=O)₂NR¹³R¹⁴, NR²⁶R²⁷, acetyl, C(=O)NR¹³R¹⁴, C(=O)OR¹³, C₁-C₆ alkyl, OCHF₂, SCF₃, OCF₃, and -C(=NH)NH₂;

~~alternatively, R²⁹ and R⁹ are taken together to form a 5-to-7 membered fused heterocyclic ring containing 1-2 heteroatom(s) each independently a member selected from the group consisting of N, O and S; wherein said 5-to-7 membered fused heterocyclic ring is substituted with 0-2 R¹⁹;~~

each R³⁰ is independently a member selected from the group consisting of H, C₃-C₇ cycloalkyl, C₁-C₄ alkyl substituted with 0-1 R²⁵, and a phenyl substituted with 0-3 R¹⁵;

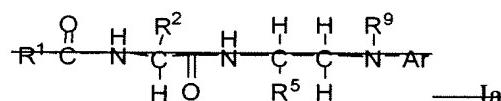
and with the proviso that R³, R⁴, R⁵, R⁶, R⁷, and R⁸, and R⁹ are not all hydrogen.

2-3. (canceled)

4. (previously presented) The compound of claim 1, wherein R¹ is phenyl substituted with 0-3 R^{1a}.

5-6. (canceled)

7. (currently amended) The compound of claim 9 ~~1~~, according to formula Ia:



wherein:

R¹ is C₃-C₈ cycloalkyl substituted with 0-2 R^{1b}, wherein said C₃-C₈ cycloalkyl is saturated or unsaturated; and

R² is a member selected from the group consisting of a phenyl substituted with 0-3 R¹⁵, a C₁-C₆ alkyl substituted with 0-2 R^{2a}, and a C₃-C₇ cycloalkyl substituted with 0-2 R¹⁹; and

~~Ar is phenyl substituted with 0-3 R²⁹, or alternatively, R²⁹ and R⁹ are taken together to form a 5-to-7 membered fused heterocyclic ring containing 1-2 heteroatom(s) each independently a member selected from the group consisting of N, O and S; wherein said 5-to-7 membered fused heterocyclic ring is substituted with 0-2 R¹⁹.~~

8. (currently amended) The compound of claim [[7]] 9, wherein:

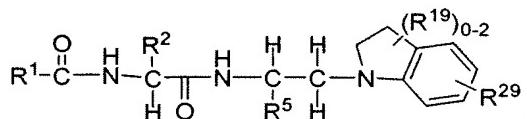
R² is a member selected from the group consisting of a C₁-C₂ alkyl substituted with 1 R^{2a}, and C₁-C₆ alkyl;

each R^{2a} is independently a member selected from the group consisting of a phenyl substituted with 0-3 R¹⁵, and a C₃-C₈ cycloalkyl substituted with 0-2 R¹⁹;

R⁵ is a member selected from the group consisting of H, C₃-C₇ cycloalkyl; and a C₁-C₆ alkyl substituted with 0-1 R¹⁸, wherein said C₁-C₆ alkyl optionally contains a heteroatom selected from the group consisting of O, S, S(=O), S(=O)₂ and NR¹⁷; and

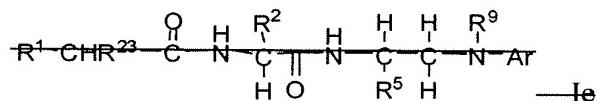
each R¹⁸ is independently a member selected from the group consisting of H, OH, F, Cl, CN, C(=O)OR³⁰, C(=O)NR¹³R¹⁴, NR¹¹R¹², a phenyl substituted with 0-3 R¹⁵, and C₃-C₈ cycloalkyl.

9. (currently amended) The compound of claim [[7]] 1, wherein said compound is of the formula:



10-15. (canceled)

16. (currently amended) The compound of claim 18 1, according to formula Ie



wherein:

R¹ is C₆-C₁₀ aryl substituted with 0-3 R^{1a}; and

each R^{1a} is independently a member selected from the group consisting of H, C₁-C₃ perfluoroalkyl, C₃-C₇ cycloalkyl, F, Cl, Br, CN, NO₂, OR¹⁰, SCH₃, S(=O)CH₃, S(=O)₂R¹⁰, NR¹¹R¹², acetyl, C(=O)OR¹³, C(=O)NR¹³R¹⁴, S(=O)₂NR¹³R¹⁴, phenyl substituted with 0-3 R¹⁵, and a C₁-C₄ alkyl substituted with 0-2 R¹⁶; and

Ar is phenyl substituted with 0-3 R²⁹, or alternatively, R²⁹ and R⁹ are taken together to form a 5-to-7-membered fused heterocyclic ring containing 1-2 heteroatom(s) each independently a member selected from the group consisting of N, O and S; wherein said 5-to-7-membered fused heterocyclic ring is substituted with 0-2 R¹⁹.

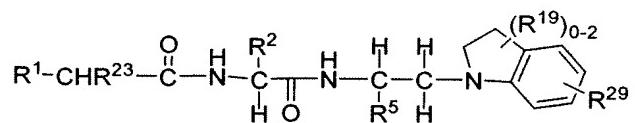
17. (currently amended) The compound of claim [[16]] 18, wherein:

R² is a member selected from the group consisting of a C₁-C₂ alkyl substituted with 1 R^{2a}, and C₁-C₆ alkyl;

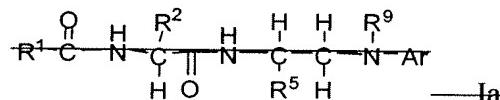
each R^{2a} is independently a member selected from the group consisting of a phenyl substituted with 0-3 R¹⁵, and a C₃-C₈ cycloalkyl substituted with 0-2 R¹⁹; and

R⁵ is a member selected from the group consisting of H, C₃-C₇ cycloalkyl; and a C₁-C₆ alkyl, wherein said C₁-C₆ alkyl optionally contains a heteroatom selected from the group consisting of O, S, S(=O), S(=O)₂ and NR¹⁷.

18. (currently amended) The compound of claim [[16]] 1, wherein said compound is of the formula:



19. (currently amended) The compound of claim 9 1, according to formula Ia



wherein:

R¹ is C₆-C₁₀ aryl substituted with 0-3 R^{1a};

each R^{1a} is independently a member selected from the group consisting of H, C₁-C₃ perfluoroalkyl, C₃-C₇ cycloalkyl, F, Cl, Br, CN, NO₂, OR¹⁰, SCH₃, S(=O)CH₃, S(=O)₂R¹⁰, NR¹¹R¹², acetyl, C(=O)OR¹³, C(=O)NR¹³R¹⁴, S(=O)₂NR¹³R¹⁴, phenyl substituted with 0-3 R¹⁵; and a C₁-C₄ alkyl substituted with 0-2 R¹⁶;

R² is a member selected from the group consisting of a phenyl substituted with 0-3 R¹⁵; a C₁-C₂ alkyl substituted [[with R^{2a}]] with 1 R^{2a}, and a C₃-C₇ cycloalkyl substituted with 0-2 R¹⁹; and

each R^{2a} is independently a member selected from the group consisting of a C₆-C₁₀ aryl substituted with 0-3 R¹⁵; a C₃-C₈ cycloalkyl substituted with 0-2 R¹⁹; and a C₇-C₁₁ bicycloalkyl substituted with 0-2 R¹⁹; and

Ar is phenyl substituted with 0-3 R²⁹, or alternatively, R²⁹ and R⁹ are taken together to form a 5- to 7-membered fused heterocyclic ring containing 1-2 heteroatom(s) each

~~independently a member selected from the group consisting of N, O and S; wherein said 5-to 7-membered fused heterocyclic ring is substituted with 0-2 R¹⁹.~~

20-26. (canceled)

27. (currently amended) A pharmaceutical composition comprising the compound of Formula I(i) or I(j) in claim 1[: or] and a pharmaceutically acceptable salt and an excipient.

28. (currently amended) A pharmaceutical composition comprising the compound of claim 38 and a pharmaceutically acceptable excipient.

29-37. (canceled)

38. (currently amended) The compound of claim 1, selected from the group consisting of:

(S)-N-{1-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-4-phenoxy-benzamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-[2-(4-methoxy-phenyl)-acetylamino]-propionamide;

(S)-N-{1-[2-(5-Chloro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(7-methoxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(6-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(7-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-cyano-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

Inventors: Liu, *et al.*

Filing Date: March 23, 2004

Response to Advisory Action dated October 5, 2007

Date: October 12, 2007

Cyclopropanecarboxylic acid (S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-amide;

(S)-N-{3-Cyclohexyl-1-[2-(4-methoxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-methoxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-benzyloxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

~~N-(1-(S)-[2-(4-Methoxy phenylamino) propylecarbamoyl] 3-methyl butyl)-3-methyl benzamide;~~

~~N-(1-(S)-[2-(4-Methoxy phenylamino)-1-methyl ethylecarbamoyl] 3-methyl butyl)-3-methyl benzamide;~~

~~N-(1-(S)-[2-(4-Methoxy phenylamino)-1-(S)-methyl ethylecarbamoyl] 3-methyl butyl)-3-methyl benzamide;~~

~~N-(1-(S)-[2-(4-Methoxy phenylamino)-1-(R)-methyl ethylecarbamoyl] 3-methyl butyl)-3-methyl benzamide;~~

~~N-(2-Cyclohexyl-(1S)-[2-(4-methoxy phenylamino)-(1R)-methyl ethylecarbamoyl]-ethyl)-3-methoxy-benzamide;~~

N-{(1S)-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-1,1-dimethyl-ethylcarbamoyl]-2-phenyl-ethyl}-3-methyl-benzamide;

~~N-(1-(S)-[1-(R)-Benzylloxymethyl]-2-(4-methoxy phenylamino)-ethylcarbamoyl]-3-methyl butyl)-3-methyl benzamide;~~

~~N-(S)-{[1-(R)-Benzylloxymethyl]-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-phenyl methyl}-3-methoxy benzamide;~~

~~N-[1-(S)-[1-(R)-Benzylloxymethyl]-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-2-(4-fluoro-phenyl)-ethyl]-3-methoxy-benzamide;~~

~~N-(1-(S)-[(2-Benzyloxy-1-(R)-(5-fluoro-2,3-dihydro-indol-1-yl)methyl)-ethylcarbamoyl]-3-cyclohexyl-propyl)-3-methoxy-benzamide;~~

N-{3-Cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(R)-hydroxymethyl-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

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N-{3-Cyclohexyl-1-(R)-[(S)-2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(R)-hydroxymethyl-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S,S)-5-(5-Fluoro-2,3-dihydro-indol-1-yl)-4-[4-methyl-2-(3-methyl-benzoylamino)-pentanoylamino]-pentanoic acid benzyl ester;

(S,S)-5-(5-Fluoro-2,3-dihydro-indol-1-yl)-4-[4-methyl-2-(3-methyl-benzoylamino)-pentanoylamino]-pentanoic acid;

(S,S)-*N*-{1-[3-Carbamoyl-1-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-propylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;

~~(S,S)-N-{1-[1-(5-Fluoro-2,3-dihydro-indol-1-ylmethyl)-3-ureido-propylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;~~

(S,S)-3-[4-Cyclohexyl-2-(3-methoxy-benzoylamino)-butyrylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid benzyl ester;

(S,S)-3-[4-Cyclohexyl-2-(3-methoxy-benzoylamino)-butyrylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid;

(S,S)-*N*-{1-[1-Benzyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;

(S,S)-*N*-{3-Cyclohexyl-1-[1-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-3-methyl-butylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S,S)-*N*-{3-Cyclohexyl-1-[1-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-2-methyl-propylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S,S)-*N*-{3-Cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-phenyl-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

~~*N*{1-(S)-[2-(R)-Benzylxy-1-(R)-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-propylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;~~

~~*N*{1-(R)-[1-(R)-Benzylsulfanyl-methyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;~~

~~(S,S)-[5-[4-Cyclohexyl-2-(3-methoxy-benzoylamino)-butyrylamino]-6-(5-fluoro-2,3-dihydro-indol-1-yl)-hexyl]-carbamic acid benzyl ester;~~

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-(2-fluoro-biphenyl-4-yl)-propionamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-p-tolyl-propionamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-o-tolyl-propionamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-(4-fluoro-phenyl)-propionamide;

2-(4-Chloro-phenyl)-N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-propionamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-(R)-phenyl-propionamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-3-methyl-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-4-(methanesulfonylamino-methyl)-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-3-methanesulfonyl-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-4-methanesulfonylamino-benzamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-(4-hydroxy-phenyl)-propionamide;

4-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(S)-(2-(R)-phenyl-propionylamino)-butyramide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-(R)-phenyl-butyramide;

~~N-{1-(S)-[1-(R)-Benzylloxymethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-2-cyclohexyl-ethyl}-3-methoxy-benzamide;~~

N-{2-Cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(R)-hydroxymethyl-ethylcarbamoyl]-ethyl}-3-methoxy-benzamide;

~~N-{1-(S)-[1-(R)-Benzylloxymethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3,3-dimethyl-butyl}-3-methoxy-benzamide;~~

N-{1-(S)-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-1-(R)-hydroxymethyl-ethylcarbamoyl]-3,3-dimethyl-butyl}-3-methoxy-benzamide;

~~3-(S)-(2-(S)-Benzyloxycarbonylamino-4,4-dimethyl-pentanoylamino)-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid tert-butyl ester;~~

~~3-(S)-(2-(S)-Benzylcarbamoyl-amino-4,4-dimethyl-pentanoylamino)-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid;~~

~~4-(5-Fluoro-2,3-dihydro-indol-1-yl)-3-(S)-[2-(S)-(3-methoxy-benzoylamino)-4,4-dimethyl-pentanoylamino]-butyric acid tert-butyl ester;~~

~~3-(S)-[3-Cyclohexyl-2-(S)-(3-methoxy-benzoylamino)-propionylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid benzyl ester;~~

~~3-(S)-[3-Cyclohexyl-2-(S)-(3-methoxy-benzoylamino)-propionylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid;~~

~~4-(5-Fluoro-2,3-dihydro-indol-1-yl)-3-(S)-[2-(S)-(3-methoxy-benzoylamino)-4,4-dimethyl-pentanoylamino]-butyric acid ethyl ester;~~

~~N-{1-(S)-[2-Cyano-1-(S)-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-ethylcarbamoyl]-3,3-dimethyl-butyl}-3-methoxy-benzamide;~~

~~N-{1-(S)-[5-Amino-1-(S)-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-pentylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;~~

~~3-(S)-(2-(S)-Benzylcarbamoyl-amino-3-cyclohexyl-propionylamino)-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid benzyl ester;~~

~~1-(S)-[1-(R)-Benzylcarbamoyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-2-cyclohexyl-ethyl carbamic acid benzyl ester;~~

~~N-{3-Cyclohexyl-1-(S)-[2-(3,5-dimethoxy-benzyl)-1-(R)-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;~~

~~4-[2-(R)-[4-Cyclohexyl-2-(S)-(3-methoxy-benzoylamino)-butyryl-amino]-3-(5-fluoro-2,3-dihydro-indol-1-yl)-propoxymethyl]-benzoic acid methyl ester;~~

~~(S,S)-N-{3-Cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(4-hydroxy-benzyl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;~~

~~{2-Cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(S)-methyl-ethylcarbamoyl]-ethyl}-carbamic acid benzyl ester;~~

~~4-Benzylcarbamoyl-N-(R,S)-{[2-(4-amidinophenylamino)-1-(S)-methyl-ethylcarbamoyl]-{(2,4-dichlorophenyl)methyl}}-benzamide;~~

~~{1-(S)-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-1-(S)-methyl-ethylcarbamoyl]-3,3-dimethyl-butyl}-carbamic acid benzyl ester;~~

~~Cyclopropanecarboxylic acid {1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(S)-methyl-ethylcarbamoyl]-3,3-dimethyl-butyl}-amide;~~

(*S,S*)-2-(3-Chloro-benzenesulfonylamino)-3-cyclohexyl-N-[1-methyl-2-(4-trifluoromethoxy-phenylamino)-ethyl]-propionamide;

(*S,S*)-3-Cyclohexyl-N-[1-methyl-2-(4-trifluoromethoxy-phenylamino)-ethyl]-2-(3-trifluoromethoxy-benzenesulfonylamino)-propionamide;

N-((*S*)-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)(cyclohexyl)methyl)-3-methylbenzamide;

N-((*S*)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-(2-chlorophenyl)ethyl)-3-methylbenzamide;

N-((*S*)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-(3-chlorophenyl)ethyl)-3-methylbenzamide;

N-((*S*)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-(4-chlorophenyl)ethyl)-3-methylbenzamide;

(*S*)-N-{2-Cyclopentyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-3-methyl-benzamide;

N-((*S*)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-3,3-dimethylbutyl)-3-methylbenzamide;

N-((*S*)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-3-cyclohexylpropyl)-3-methylbenzamide;

N-((*S*)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-phenylethyl)-3-methylbenzamide;

N-(*R,S*)-((3-(5-fluoroindolin-1-yl)-1-hydroxypropan-2-(*R*)-ylcarbamoyl)(2,4-dichlorophenyl)methyl)-3,4-difluorobenzamide;

N-*(S)* ((3-(benzyloxy)-1-(5-fluoroindolin-1-yl)propan-2-(*R*)-ylcarbamoyl)(2,4-dichlorophenyl)methyl)-3,4-difluorobenzamide;

(*R,S*)-N-((2-(5-fluoroindolin-1-yl)ethylcarbamoyl)(2,4-dichlorophenyl)methyl)-3-methylbenzamide;

(*S,S*)-N-((3-(5-fluoroindolin-1-yl)-1-hydroxypropan-2-ylcarbamoyl)(2,4-dichlorophenyl)methyl)-3,4-difluorobenzamide;

(*S,S*)-4-(5-Fluoro-2,3-dihydro-indol-1-yl)-3-[2-(3-methoxy-benzoylamino)-4,4-dimethyl-pentanoylamino]-butyric acid; and

(*S*)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(5-isoxazol-3-yl-thiophene-2-sulfonylamino)-propionamide;

(S)-2-(3-Biphenyl-4-yl-ureido)-3-cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-propionamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-phenoxy-benzenesulfonylamino)-propionamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(naphthalene-1-sulfonylamino)-propionamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-trifluoromethyl-benzenesulfonylamino)-propionamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-trifluoromethoxy-benzenesulfonylamino)-propionamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-[4-(4-fluoro-phenoxy)-benzenesulfonylamino]-propionamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4'-methoxy-biphenyl-4-sulfonylamino)-propionamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-methoxy-benzenesulfonylamino)-propionamide;

(S)-3-Cyclohexyl-2-(4-difluoromethoxy-benzenesulfonylamino)-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-propionamide; ,

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-phenylmethanesulfonylamino-propionamide; ,

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(toluene-3-sulfonylamino)-propionamide; ,

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-[4-(4-methoxy-phenoxy)-benzenesulfonylamino]-propionamide; ,

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(3-methoxy-benzenesulfonylamino)-propionamide; ,

(S,S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-2-(toluene-3-sulfonylamino)-propionamide; ,

(S,S)-3-[4-(4-Dimethyl-2-(toluene-3-sulfonylamino)-pentanoylamino)-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid-tert-butyl ester]; ,

(S,S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-2-(3-trifluoromethoxy-benzenesulfonylamino)-propionamide; ,

(S,S)-2-(3-Chloro-benzenesulfonylamino)-3-cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-propionamide;

(S,S)-N-{3-Cyclohexyl-1-[1-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-3-hydroxy-propylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S,S)-3-[4,4-Dimethyl-2-(toluene-3-sulfonylamino)pentanoylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)butyric acid;

(S,S)-2-Benzenesulfonylamino-3-cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-propionamide; and

(S,S)-4,4-Dimethyl-2-(toluene-3-sulfonylamino)pentanoic acid [2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-amide.